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# A Dimensionality Reduction Approach for Qualitative Preference Aggregation

Quentin Brabant<sup>1</sup>, Miguel Couceiro<sup>1</sup>, Fabien Labernia<sup>2</sup>, Amedeo Napoli<sup>1</sup>

<sup>1</sup> LORIA (CNRS, Inria Nancy Grand Est - Université de Lorraine  
{quentin.brabant, miguel.couceiro, amedeo.napoli}@inria.fr

<sup>2</sup> LAMSADE (CNRS - Université Paris-Dauphine)  
fabien.labernia@dauphine.fr

## 1 Qualitative preference aggregation models

In this paper we briefly present a method for reducing the dimensionality of data in a qualitative preference aggregation framework. For a more complete description of this approach, see [4]. For an alternative approach based on rough sets theory, see [1].

We consider the following setting.  $X$  is a set of alternatives that are evaluated according to a set of criteria represented by their indices:  $[n] = \{1, \dots, n\}$ . For an alternative  $x \in X$  we denote by  $(x_1, \dots, x_n) \in L^n$  the tuple of the evaluations of  $x$  in each criterion.  $L$  is called the *evaluation space*, and is a distributive lattice for which we denote respectively by 0 and 1 the minimal and maximal element. We consider a binary preference relation  $\preceq$  between the alternatives that can be expressed in terms of a *utility function*:

$$\forall x, y \in X : x \preceq y \Rightarrow U(x) \leq U(y),$$

where  $U : X \rightarrow L$  associates a global evaluation on  $L$  to each alternative, and is obtained through the aggregation of the evaluations in criteria by a Sugeno integral  $\mathcal{S}_\mu : L^n \rightarrow L$ . In other words we have  $U(x) = \mathcal{S}_\mu(x_1, \dots, x_n)$ . The Sugeno integral defined over distributive lattices [3], is expressed

$$\mathcal{S}_\mu(x_1, \dots, x_n) = \bigvee_{I \subseteq [n]} \mu(I) \bigwedge_{i \in I} x_i,$$

where  $\mu : 2^{[n]} \rightarrow L$  a capacity, that is to say a non-decreasing set function on  $[n]$ , with  $\mu(\emptyset) = 0$  and  $\mu([n]) = 1$ . Capacities (and Sugeno integrals) are defined by a value on  $L$  for each subset of  $[n]$ , and therefore carry an intrinsic complexity, that grows exponentially with  $n$ . We now consider a set  $\mathcal{D} = \{(\mathbf{x}^1, y^1), \dots, (\mathbf{x}^m, y^m)\} \subseteq L^n \times L$ , where each  $\mathbf{x}^i = (x_1^i, \dots, x_n^i) \in L^n$  is a tuple of evaluations in  $n$  criteria, and  $y^i$  is a utility value associated to  $\mathbf{x}^i$ . We want to learn a Sugeno integral  $\mathcal{S}_\mu$  that generalizes these data. Ideally this function would be such that  $\mathcal{S}_\mu(\mathbf{x}^j) = y^j$  for any  $j \in \{1, \dots, m\}$ . However, it is very common that no such function exists: in that case  $\mathcal{D}$  is said to be *inconsistent*, and we aim at learning a Sugeno integral that realizes the prediction of  $y^j$  for each element, with an error as low as possible. Because of the nature of capacities, this optimization problem is on  $2^n$  variables, and is therefore hard to solve when a high number of criteria is considered.

## 2 Dimensionnality reduction based on quality measure

By a *quality measure* over  $\mathcal{D}$  we mean a degree with which  $\mathcal{D}$  satisfies a certain hypothesis. In this presentation we consider two of such measures.

The first quality measure is the *monotonicity degree*, that is, the ratio of pairs  $\{i, j\} \subseteq \{1, \dots, m\}$  that satisfy the following condition

$$y^i > y^j \Rightarrow \exists k \in [n] : y_k^i > y_k^j.$$

This condition can be seen as a generalization of the Pareto condition to partially ordered evaluation spaces. The second quality measure is the *compatibility degree*, that is, the ratio of pairs satisfying the condition

$$\exists \mathcal{S}_\mu : [\mathcal{S}_\mu(\mathbf{x}^i) = y^i \text{ and } \mathcal{S}_\mu(\mathbf{x}^j) = y^j]. \quad (1)$$

This condition is justified by results from [2] that apply only when  $L$  is totally ordered. Indeed it can be shown that  $\mathcal{D}$  is consistent if and only if (2) is true for any pair from  $\mathcal{D}$ . Moreover, for a given pair this condition can be checked in a linear time w.r.t.  $n$ . Hence, provided that  $L$  is totally ordered, the compatibility degree is both theoretically meaningful and practically interesting. If  $L$  is not totally ordered, the monotonicity degree is the quality measure that makes sense.

The principle of the algorithm for dimensionality reduction that we propose is to iteratively remove a criterion, in order to minimize the decrease of the quality of the dataset at each step. Criteria are deleted until it is impossible to remove a criterion without decreasing the quality of the data below a certain ratio  $\alpha$ . This algorithm was tested on empirical data<sup>1</sup> and allowed a reduction of the number of criteria from 7 to 3. Aggregation models trained on original data and on data reduced to 3 criteria showed to have similar accuracy. On the other hand, models trained on data with only 2 criteria left had significantly worse accuracy, suggesting that a reduction to 3 criteria constitutes the best compromise between simplicity and accuracy for these data.

Future research work should include further empirical studies and should aim to determining a procedure for deciding the optimal value of  $\alpha$ , currently being set by hand.

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<sup>1</sup>Tripadvisor: <http://sifaka.cs.uiuc.edu/~wang296/Data/index.html>